Amendments to the Claims

The following listing of claims will replace all prior versions and listings of claims in the application.

Listing of Claims:

1. (Currently amended) A compound of formula (I):

$$(R^3)_{m}$$
 R^6 R^6 R^5

wherein:

m is an integer from 0 to 4;

 R^1 is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-OR^7$ or $-N(R^8)R^9$, with the proviso that R^1 is not 3- or 4- (1,1,1,3,3,3-hexafluoro-2-hydroxy-2-propyl)phenyl;

R², R⁴, R⁵ and R⁶ are selected from (a) and (b) as follows:

(a) R² and R⁶ are selected from (i) and (ii) as follows: (i) R² and R⁶ are each independently hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl; or (ii) R² and R⁶ together form optionally substituted alkylene or optionally substituted alkenylene; and

R⁴ and R⁵ are selected from (i) and (ii) as follows: (i) R⁴ and R⁵ are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted

alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-N(R^8)R^9$, $-OR^7$, $-S(O)_jR^{11}$ where j is 1 or 2, $-B(R^{22})_2$, $-P(R^{22})_2$, $-P(O)(R^{22})_2$ and $-C(E)R^{23}$, where E is selected from O, S and NR^7 ; or (ii) R^4 and R^5 together form optionally substituted alkylene, optionally substituted alkylene or optionally substituted alkyleneazaalkylene; or

(b) R² and R⁵, or R² and R⁴, or R⁶ and R⁵, or R⁶ and R⁴, together form a 4, 5, 6 or 7 membered optionally substituted heterocyclyl group, or a 5 or 6 membered optionally substituted heteroaryl group; and the remainder of R², R⁴, R⁵ and R⁶ are each independently selected as in (i) above;

each R³ is independently selected from the group consisting of halo, pseudohalo, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, - N(R¹²)R¹³, -OR¹⁴, -C(E)R¹⁵ where E is O, S or NR⁷, and -S(O)_vR¹⁶ where y is 0, 1 or 2;

or any two R³ groups, which substitute adjacent carbons on the ring, together form optionally substituted alkylene, optionally substituted alkenylene, optionally substituted alkylenedioxy, optionally substituted thioalkylenoxy, or optionally substituted alkylenedithioxy;

each R⁷ is independently selected from the group consisting of hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted recordly, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, and optionally substituted heterocyclylalkyl;

R⁸ and R⁹ are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted aryl,

optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-S(O)_jR^{10}$ where j is 1 or 2, and $-C(M)R^{11}$, where M is selected from O and S;

or R⁸ and R⁹ together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene;

each R¹⁰ is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, and optionally substituted heterocyclylalkyl;

each R^{11} is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkyll, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-OR^{10}$ and $-N(R^7)_2$;

 R^{12} and R^{13} are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -C(M) R^{17} where M is O or S, and -S(O) $_iR^{18}$ where j is 1 or 2;

or R¹² and R¹³ together form optionally substituted alkylene, optionally substituted alkenylene, optionally substituted alkyleneoxyalkylene or optionally substituted alkyleneazaalkylene;

R¹⁴ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted

aralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl or -C(M)R¹⁷ where M is O or S;

R¹⁵ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OH, -OR¹⁴ or -N(R¹²)R¹³;

R¹⁶ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OH, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁷ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, oPR¹⁹ or -N(R²⁰)R²¹;

 R^{18} is optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-OR^{19}$ or $-N(R^{20})R^{21}$;

R¹⁹ is alkyl, alkenyl, alkynyl, cycloalkyl, heterocyclyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl or heteroaralkyl;

 R^{20} and R^{21} are each independently selected from hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, heterocyclyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl and heteroaralkyl,

or R^{20} and R^{21} together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene;

each R^{22} is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-OR^7$ and $-N(R^7)_2$;

 R^{23} is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-OR^{10}$, $-N(R^7)_2$, or $-N(R^7)N(R^7)_2$;

wherein each of the above R¹-R²³ groups, when substituted, are substituted with one or more substituents each independently selected from Q^1 , where Q^1 is halo, pseudohalo, hydroxy, oxo, thia, nitrile, nitro, formyl, mercapto, carboxy, carboxyalkyl, alkyl, haloalkyl, polyhaloalkyl, aminoalkyl, diaminoalkyl, alkenyl containing 1 to 2 double bonds, alkynyl containing 1 to 2 triple bonds, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl, aralkynyl, aralkynyl, heteroarylalkyl, trialkylsilyl, dialkylarylsilyl, alkyldiarylsilyl, triarylsilyl, alkylidene, arylalkylidene, alkylcarbonyl, cycloalkylcarbonyl, heterocyclylcarbonyl, arylcarbonyl, heteroarylcarbonyl, alkoxycarbonyl, alkoxycarbonylalkyl, aryloxycarbonyl, aryloxycarbonylalkyl, aralkoxycarbonyl, aralkoxycarbonylalkyl, arylcarbonylalkyl, aminocarbonyl, alkylaminocarbonyl, dialkylaminocarbonyl, arylaminocarbonyl, diarylaminocarbonyl, arylalkylaminocarbonyl, alkoxy, aryloxy, heteroaryloxy, heteroaralkoxy, heterocyclyloxy, heterocyclylalkoxy, cycloalkoxy, perfluoroalkoxy, alkenyloxy, alkynyloxy, aralkoxy, alkylcarbonyloxy, arylcarbonyloxy, aralkylcarbonyloxy, alkoxycarbonyloxy, aryloxycarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, alkylarylaminocarbonyloxy, diarylaminocarbonyloxy, guanidino, isothioureido, amidino, alkylamidino, arylamidino, aminothiocarbonyl, alkylaminothiocarbonyl, arylaminothiocarbonyl, amino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, arylaminoalkyl, diarylaminoalkyl,

alkylarylamino, dialkylamino, dialkylamino, haloalkylamino, arylamino, diarylamino, alkylarylamino, alkylcarbonylamino, alkoxycarbonylamino, aralkoxycarbonylamino, arylcarbonylamino, arylcarbonylaminoalkyl, aryloxycarbonylaminoalkyl, aryloxyarylcarbonylamino, aryloxycarbonylamino, alkylsulfonylamino, arylsulfonylamino, $heteroarylsulfonylamino,\ heteroarylthio,\ azido,\ -N^+(R^{24})_3,\ -P(R^{25})_2,$ $-P(O)(R^{25})_2$, $-OP(O)(R^{25})_2$, $-N(R^{24})C(O)R^{26}$, dialkylphosphonyl, alkylarylphosphonyl, diarylphosphonyl, hydroxyphosphonyl, alkylthio, arylthio, perfluoroalkylthio, carboxyalkylthio, thiocyano, isothiocyano, alkylsulfinyloxy, alkylsulfonyloxy, arylsulfinyloxy, arylsulfinyloxy, hydroxysulfonyloxy, alkoxysulfonyloxy, aminosulfonyloxy, alkylaminosulfonyloxy, dialkylaminosulfonyloxy, arylaminosulfonyloxy, diarylaminosulfonyloxy, alkylarylaminosulfonyloxy, alkylsulfinyl, alkylsulfonyl, arylsulfinyl, arylsulfonyl, hydroxysulfonyl, alkoxysulfonyl, aminosulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, arylaminosulfonyl, diarylaminosulfonyl or alkylarylaminosulfonyl; or two O¹ groups, which substitute atoms in a 1,2 or 1,3 arrangement, together form alkylenedioxy (i.e., -O-(CH₂)_v-O-), thioalkylenoxy (i.e., _-S-(CH₂)_y-O-), or alkylenedithioxy (i.e., -S-(CH₂)_y-S-), where y is 1 or 2; or two O¹ groups, which substitute the same atom, together form alkylene;

each R²⁴ is independently selected from the group consisting of hydrogen, alkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, heterocyclyl and heterocyclylalkyl;

each R²⁵ is independently selected from the group consisting of hydroxy, alkoxy, aralkoxy, alkyl, heterocyclyl, aryl and -N(R²⁷)R²⁸,

 R^{26} is alkoxy, aralkoxy, alkyl, heteroaryl, heterocyclyl, aryl or -N(R^{27}) R^{28} ; R^{27} and R^{28} are each independently hydrogen, alkyl, aralkyl, aryl, heteroaryl, heteroaralkyl or heterocyclyl,

or R²⁷ and R²⁸ together form alkylene, azaalkylene, oxaalkylene or thiaalkylene; and each Q¹ is optionally substituted by one or more substituents selected from Q²; where each Q² is independently halo, pseudohalo, hydroxy, oxo, thia, nitrile, nitro, formyl, mercapto, carboxy, carboxyalkyl, alkyl, haloalkyl, polyhaloalkyl, aminoalkyl, diaminoalkyl, alkenyl containing 1 to 2 double bonds, alkynyl containing 1 to 2 triple bonds, cycloalkyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl, aralkenyl, aralkynyl,

heteroarylalkyl, trialkylsilyl, dialkylarylsilyl, alkyldiarylsilyl, triarylsilyl, alkylidene, arylalkylidene, alkylcarbonyl, arylcarbonyl, heteroarylcarbonyl, alkoxycarbonyl, alkoxycarbonylalkyl, aryloxycarbonyl, aryloxycarbonylalkyl, aralkoxycarbonyl, aralkoxycarbonylalkyl, arylcarbonylalkyl, aminocarbonyl, alkylaminocarbonyl, dialkylaminocarbonyl, arylaminocarbonyl, diarylaminocarbonyl, arylalkylaminocarbonyl, alkoxy, aryloxy, heteroaryloxy, heteroaralkoxy, heterocyclyloxy, heterocyclylalkoxy, cycloalkoxy, perfluoroalkoxy, alkenyloxy, alkynyloxy, aralkoxy, alkylcarbonyloxy, arylcarbonyloxy, aralkylcarbonyloxy, alkoxycarbonyloxy, aryloxycarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, alkylarylaminocarbonyloxy, diarylaminocarbonyloxy, guanidino, isothioureido, amidino, alkylamidino, arylamidino, aminothiocarbonyl, alkylaminothiocarbonyl, arylaminothiocarbonyl, amino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, arylaminoalkyl, diarylaminoalkyl, alkylarylamino, dialkylamino, dialkylamino, haloalkylamino, arylamino, diarylamino, alkylarylamino, alkylcarbonylamino, alkoxycarbonylamino, aralkoxycarbonylamino, arylcarbonylamino, arylcarbonylaminoalkyl, aryloxycarbonylaminoalkyl, aryloxyarylcarbonylamino, aryloxycarbonylamino, alkylsulfonylamino, arylsulfonylamino, heteroarylsulfonylamino, heterocyclylsulfonylamino, heteroarylthio, azido, -N⁺(R²⁴)₃, -P(R²⁵)₂, $-P(O)(R^{25})_2$, $-OP(O)(R^{25})_2$, $-N(R^{24})C(O)R^{26}$, dialkylphosphonyl, alkylarylphosphonyl, diarylphosphonyl, hydroxyphosphonyl, alkylthio, arylthio, perfluoroalkylthio, carboxyalkylthio, thiocyano, isothiocyano, alkylsulfinyloxy, alkylsulfonyloxy, arylsulfinyloxy, arylsulfinyloxy, hydroxysulfonyloxy, alkoxysulfonyloxy, aminosulfonyloxy, alkylaminosulfonyloxy, dialkylaminosulfonyloxy, arylaminosulfonyloxy, diarylaminosulfonyloxy, alkylarylaminosulfonyloxy, alkylsulfinyl, alkylsulfonyl, arylsulfinyl, arylsulfonyl, hydroxysulfonyl, alkoxysulfonyl, aminosulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, arylaminosulfonyl, diarylaminosulfonyl or alkylarylaminosulfonyl; or two Q² groups, which substitute atoms in a 1,2 or 1,3 arrangement, together form alkylenedioxy (i.e., -O-(CH₂)_v-O-), thioalkylenoxy (i.e., -S-(CH₂)_v-O-) or alkylenedithioxy (i.e., -S-(CH₂)_v-S-) where y is 1 or 2; or two Q² groups, which substitute the same atom, together form alkylene,

where R²⁴, R²⁵, R²⁶, R²⁷ and R²⁸ are as defined above;

as a stereoisomer, racemate or mixture thereof; or as a pharmaceutically acceptable derivative salt thereof;

with the proviso that the compound of formula (I) can not be N-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide.

2. (Currently amended) The compound of claim 1 wherein the compound of formula (I) is a A compound of formula (II):

$$(R^3)_{m}$$
 $(R^{1a})_{n}$ $(R^{1a})_{n}$ $(R^{1a})_{n}$ $(R^{1a})_{n}$

wherein:

n is an integer from 0 to 5;

m is an integer from 0 to 4;

each R^{1a} and R^3 are independently selected from the group consisting of halo, pseudohalo, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-N(R^{12})R^{13}$, $-OR^{14}$, $-C(E)R^{15}$ where E is O, S or NR^7 , and $-S(O)_yR^{16}$ where y is 0, 1 or 2, with the proviso that R^{1a} is not 3- or 4-C(OH)(CF₃)₂;

or any two R^{1a} groups or R³ groups, which substitute adjacent carbons on the ring, together with atoms to which they are attached, form optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substitute

R², R⁴, R⁵ and R⁶ are selected from (a) and (b) as follows:

(a) R² and R⁶ are selected from (i) and (ii) as follows: (i) R² and R⁶ are each independently hydrogen or optionally substituted alkyl; or (ii) R² and R⁶ together form alkylene or alkenylene;

 R^4 and R^5 are selected from (i) and (ii) as follows: (i) R^4 and R^5 are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-N(R^8)R^9$, $-OR^7$, $S(O)_jR^{11}$ where j is 1 or 2, $-B(R^{22})_2$, $-P(R^{22})_2$, $-P(O)(R^{22})_2$, and $-C(E)R^{23}$ where E is selected from O, S and NR^7 ; or (ii) R^4 and R^5 together form optionally substituted alkylene, optionally substituted alkylene or optionally substituted alkyleneazaalkylene;

(b) R² and R⁵, or R² and R⁴, or R⁶ and R⁵, or R⁶ and R⁴, together form a 5, 6 or 7 membered optionally substituted heterocyclyl group, or a 5 or 6 membered optionally substituted heteroaryl group; and the remainder of R², R⁴, R⁵ and R⁶ are each independently selected as in (i) above;

each R⁷ is independently selected from the group consisting of hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted recordly, optionally substituted eycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, and optionally substituted heterocyclylalkyl;

 R^8 and R^9 are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-S(O)_jR^{10}$ where j is 1 or 2, and $-C(M)R^{11}$, where M is selected from O and S;

or R⁸ and R⁹ together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene;

each R¹⁰ is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl;

each R^{11} is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-OR^{10}$ and $-N(R^7)_2$;

 R^{12} and R^{13} are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -C(M) R^{17} where M is O or S, and -S(O) $_jR^{14}$ where j is 1 or 2;

or \mathbb{R}^{12} and \mathbb{R}^{13} together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene;

R¹⁴ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl or -C(M)R¹⁷ where m is O or S;

R¹⁵ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OH, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁶ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OH, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁷ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁸ is optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted aralkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted beteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁹ is alkyl, alkenyl, alkynyl, cycloalkyl, heterocyclyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl or heteroaralkyl;

R²⁰ and R²¹ are each independently selected from hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, heterocyclyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl and heteroaralkyl,

or R²⁰ and R²¹ together form alkylene, alkenylene or alkyleneoxyalkylene; each R²² is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted eycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, -OR⁷ and -N(R⁷)₂;

R²³ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl,

optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-OR^{10}$, $-N(R^7)_2$, $-N(R^7)N(R^7)_2$;

wherein each of the above R¹-R²³ groups, when substituted, are substituted with one or more substituents each independently selected from O¹, where O¹ is halo, pseudohalo, hydroxy, oxo, thia, nitrile, nitro, formyl, mercapto, carboxy, carboxyalkyl, alkyl, haloalkyl, polyhaloalkyl, aminoalkyl, diaminoalkyl, alkenyl containing 1 to 2 double bonds, alkynyl containing 1 to 2 triple bonds, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl, aralkynyl, aralkynyl, heteroarylalkyl, trialkylsilyl, dialkylarylsilyl, alkyldiarylsilyl, triarylsilyl, alkylidene, arylalkylidene, alkylcarbonyl, arylcarbonyl, heteroarylcarbonyl, alkoxycarbonyl, alkoxycarbonylalkyl, aryloxycarbonyl, aryloxycarbonylalkyl, aralkoxycarbonyl, aralkoxycarbonylalkyl, arylcarbonylalkyl, aminocarbonyl, alkylaminocarbonyl, dialkylaminocarbonyl, arylaminocarbonyl, diarylaminocarbonyl, arylalkylaminocarbonyl, alkoxy, aryloxy, heteroaryloxy, heteroaralkoxy, heterocyclyloxy, heterocyclylalkoxy, cycloalkoxy, perfluoroalkoxy, alkenyloxy, alkynyloxy, aralkoxy, alkylcarbonyloxy, arylcarbonyloxy, aralkylcarbonyloxy, alkoxycarbonyloxy, aryloxycarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, alkylarylaminocarbonyloxy, diarylaminocarbonyloxy, guanidino, isothioureido, amidino, alkylamidino, arylamidino, aminothiocarbonyl, alkylaminothiocarbonyl, arylaminothiocarbonyl, amino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, arylaminoalkyl, diarylaminoalkyl, alkylarylaminoalkyl, alkylamino, dialkylamino, haloalkylamino, arylamino, diarylamino, alkylarylamino, alkylcarbonylamino, alkoxycarbonylamino, aralkoxycarbonylamino, arylcarbonylamino, arylcarbonylaminoalkyl, aryloxycarbonylaminoalkyl, aryloxyarylcarbonylamino, aryloxycarbonylamino, alkylsulfonylamino, arylsulfonylamino, heteroarylsulfonylamino, heterocyclylsulfonylamino, heteroarylthio, azido, $-N^+(R^{24})_3$, $-P(R^{25})_2$, $-P(O)(R^{25})_2$, $-OP(O)(R^{25})_2$, $-N(R^{24})C(O)R^{26}$, dialkylphosphonyl, alkylarylphosphonyl, diarylphosphonyl, hydroxyphosphonyl, alkylthio, arylthio, perfluoroalkylthio, carboxyalkylthio, thiocyano, isothiocyano, alkylsulfinyloxy, alkylsulfonyloxy, arylsulfinyloxy, arylsulfonyloxy, hydroxysulfonyloxy, alkoxysulfonyloxy,

aminosulfonyloxy, alkylaminosulfonyloxy, dialkylaminosulfonyloxy, arylaminosulfonyloxy, diarylaminosulfonyloxy, alkylaufinyl, alkylsulfinyl, arylsulfinyl, arylsulfonyl, hydroxysulfonyl, alkoxysulfonyl, aminosulfonyl, alkylaminosulfonyl, diarylaminosulfonyl, or alkylaminosulfonyl; or two Q^1 groups, which substitute atoms in a 1,2 or 1,3 arrangement, together form alkylenedioxy (*i.e.*, -O-(CH₂)_y-O-), thioalkylenoxy (*i.e.*, -S-(CH₂)_y-O-) or alkylenedithioxy (*i.e.*, -S-(CH₂)_y-S-) where y is 1 or 2; or two Q^1 groups, which substitute the same atom, together form alkylene;

each R²⁴ is independently selected from the group consisting of hydrogen, alkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, heterocyclyl and heterocyclylalkyl;

each R^{25} is independently selected from the group consisting of hydroxy, alkoxy, aralkoxy, alkyl, heterocyclyl, aryl and $-N(R^{27})R^{28}$,

 R^{26} is alkoxy, aralkoxy, alkyl, heteroaryl, heterocyclyl, aryl or -N(R^{27}) R^{28} ; R^{27} and R^{28} are each independently hydrogen, alkyl, aralkyl, aryl, heteroaryl, heteroaralkyl or heterocyclyl,

or R²⁷ and R²⁸ together form alkylene, azaalkylene or thiaalkylene; and each Q¹ is optionally substituted by one or more substituents selected from Q²; where each Q² is independently halo, pseudohalo, hydroxy, oxo, thia, nitrile, nitro, formyl, mercapto, carboxy, carboxyalkyl, alkyl, haloalkyl, polyhaloalkyl, aminoalkyl, diaminoalkyl, alkenyl containing 1 to 2 double bonds, alkynyl containing 1 to 2 triple bonds, cycloalkyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl, aralkenyl, aralkynyl, heteroarylalkyl, trialkylsilyl, dialkylarylsilyl, alkyldiarylsilyl, triarylsilyl, alkylidene, arylalkylidene, alkylcarbonyl, arylcarbonyl, heteroarylcarbonyl, alkoxycarbonyl, alkoxycarbonyl, aralkoxycarbonylalkyl, aryloxycarbonylalkyl, aminocarbonyl, alkylaminocarbonyl, dialkylaminocarbonyl, arylalkylaminocarbonyl, alkoxy, aryloxy, heteroaryloxy, heteroaralkoxy, heterocyclyloxy, heterocyclylalkoxy, cycloalkoxy, perfluoroalkoxy, alkenyloxy, alkynyloxy, aralkoxy, alkylcarbonyloxy, arylcarbonyloxy, aralkoxycarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy,

alkylarylaminocarbonyloxy, diarylaminocarbonyloxy, guanidino, isothioureido, amidino, alkylamidino, arylamidino, aminothiocarbonyl, alkylaminothiocarbonyl, arylaminothiocarbonyl, amino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, arylaminoalkyl, diarylaminoalkyl, alkylarylamino, dialkylamino, dialkylamino, haloalkylamino, arylamino, diarylamino, alkylarylamino, alkylcarbonylamino, alkoxycarbonylamino, aralkoxycarbonylamino, arylcarbonylamino, arylcarbonylaminoalkyl, aryloxycarbonylaminoalkyl, aryloxyarylcarbonylamino, aryloxycarbonylamino, alkylsulfonylamino, arylsulfonylamino, heteroarylsulfonylamino, heterocyclylsulfonylamino, heteroarylthio, azido, -N⁺(R²⁴)₃, -P(R²⁵)₂, $-P(O)(R^{25})_2$, $-OP(O)(R^{25})_2$, $-N(R^{24})C(O)R^{26}$, dialkylphosphonyl, alkylarylphosphonyl, diarylphosphonyl, hydroxyphosphonyl, alkylthio, arylthio, perfluoroalkylthio, carboxyalkylthio, thiocyano, isothiocyano, alkylsulfinyloxy, alkylsulfonyloxy, arylsulfinyloxy, arylsulfinyloxy, hydroxysulfonyloxy, alkoxysulfonyloxy, aminosulfonyloxy, alkylaminosulfonyloxy, dialkylaminosulfonyloxy, arylaminosulfonyloxy, diarylaminosulfonyloxy, alkylarylaminosulfonyloxy, alkylsulfinyl, alkylsulfonyl, arylsulfinyl, arylsulfonyl, hydroxysulfonyl, alkoxysulfonyl, aminosulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, arylaminosulfonyl, diarylaminosulfonyl or alkylarylaminosulfonyl; or two Q² groups, which substitute atoms in a 1,2 or 1,3 arrangement, together form alkylenedioxy (i.e., -O-(CH₂)_v-O-), thioalkylenoxy (i.e., -S-(CH₂)_v-O-) or alkylenedithioxy (i.e., -S-(CH₂)_v-S-) where y is 1 or 2; or two Q² groups, which substitute the same atom, together form alkylene;

where R²⁴, R²⁵, R²⁶, R²⁷ and R²⁸ are as defined above;

as a stereoisomer, racemate or mixture thereof.; or as a pharmaceutically acceptable derivative salt thereof.

3. (Currently amended) The compound of claim 2 wherein the compound of formula (II) is a compound of formula (III):

$$(R^3)_m$$
 $(R^{1a})_n$ $(R^{1a})_n$ $(R^{5a})_t$

wherein:

n is an integer from 0 to 5;

m is an integer from 0 to 4;

t is an integer from 0 to 5;

each R^{1a} is independently selected from the group consisting of alkyl, hydroxy, alkoxy, alkoxyalkoxy, aralkoxy, amino, alkylamino, dialkylamino, halo, haloalkyl, haloalkoxy, cyano, carboxy, alkoxycarbonyl, alkoxycarbonylalkoxy, heteroaryl, heterocyclyl, heterocyclylalkoxy, and aryl (optionally substituted by one or more substituents selected from the group consisting of alkyl, alkoxy, halo, cyano, carboxy, cyano, and alkoxycarbonyl);

R² is hydrogen or alkyl; alkyl'

each R³ is independently selected from the group consisting of alkyl, alkoxy, halo, hydroxy, aralkoxy, alkoxycarbonylalkoxy, aryl (optionally substituted by one or more substituents independently selected from the group consisting of alkyl, halo, alkoxy, carboxy, alkoxycarbonyl, cyano), heteroaryl and heterocyclyl;

R⁴ is hydrogen or alkyl;

each R^{5a} is independently selected from the group consisting of alkyl, alkoxy, halo, alkylcarbonyl, haloalkyl, haloalkoxy, aryl, cyano, carboxy, alkoxycarbonyl, nitro, and $-N(R^{24})C(O)R^{26}$:

or two adjacent R^{5a} groups form phenyl, 5-6 membered heteroaryl, O-(CH₂)_y-O-, -S-(CH₂)_y-O-, or -S-(CH₂)_y-S- an aryl, heterocyclyl or heteroaryl; and R^{6} is hydrogen or alkyl.

- 4. (Currently amended) The compound of claim 3 wherein m is 0 or 1, n is 1 and R^{1a} is independently selected from alkoxy, halo, haloalkyl, haloalkoxy, cyano, optionally substituted aryl, aryl (optionally substituted by one or more substituents independently selected from the group consisting of alkyl, halo, alkoxy, carboxy, alkoxycarbonyl, cyano), heterocyclyl, and heteroaryl.
- 5. (Currently amended) The compound of claim 4 selected from the group consisting of the following:

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzenesulfonamide; *N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-4-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzenesulfonamide;

4-methoxy-*N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzenesulfonamide;

4-chloro-*N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,*N*-dimethylbenzenesulfonamide;

4-({1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-methyl-sulfamoyl)-benzoic acid;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,*N*-dimethylbenzenesulfonamide;

4-methoxy-*N*-{1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-butyl-*N*-{1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-4-trifluoromethylbenzenesulfonamide;

2,4,6-trichloro-*N*-{1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl} methylamide;

N-{1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,*N*-dimethylbenzenesulfonamide;

4-methoxy-*N*-{1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-butyl-*N*-{1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-4-trifluoromethylbenzenesulfonamide;

2,4,6-trichloro-*N*-{1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,*N*-dimethylbenzenesulfonamide;

4-methoxy-*N*-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-butyl-*N*-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methyl-4-trifluoromethylbenzenesulfonamide;

2,4,6-trichloro-*N*-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-methoxy-*N*-methylbenzenesulfonamide;

N-{1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,*N*-dimethylbenzenesulfonamide;

N-{1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methyl-4-trifluoromethylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-methoxy-N-methylbenzenesulfonamide;

 ${\it N-\{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl\}-N-methylbenzenesulfonamide;}$

N-{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,N-dimethylbenzenesulfonamide;

 $\label{eq:N-local-state} N-\{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl\}-N-methyl-4-trifluoromethylbenzenesulfonamide;$

2,4,6-trichloro-*N*-{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,*N*-dimethylbenzenesulfonamide;

4-methoxy-*N*-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-tert-butyl-N-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methylbenzamide;

4-butyl-*N*-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methyl-4-trifluoromethylbenzenesulfonamide;

2,4,6-trichloro-*N*-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[8-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[8-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-

dihydroquinazolin-2-yl]ethyl}-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[8-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[5-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[5-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[5-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[6-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[6-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[6-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(3-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[3-(3-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(3-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(2-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-isopropyl-*N*-{1-[3-(2-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(2-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[8-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[8-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[8-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl} methylamide;

4-*tert*-butyl-*N*-{1-[3-(4-ethoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-isopropoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-isobutoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-*n*-butoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-methoxy-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide

4-*tert*-butyl-*N*-{1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-methoxy-N-methylbenzenesulfonamide;

N-{1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-tert-butyl-N-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-cyanophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,N-dimethylbenzenesulfonamide;

N-{1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-chloro-N-methylbenzenesulfonamide;

N-{1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-4-trifluoromethylbenzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-trifluoromethoxyphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-benzenesulfonamide;

4-isopropyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-trifluoromethoxyphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-benzenesulfonamide;

biphenyl-4-sulfonic acid methyl-{1-[4-oxo-3-(4-trifluoromethoxyphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-amide;

4-*tert*-butyl-*N*-{1-[3-(4-fluorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-fluorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-fluorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-trifluoromethylphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-benzenesulfonamide;

4-isopropyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-trifluoromethylphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-benzenesulfonamide;

biphenyl-4-sulfonic acid methyl-{1-[4-oxo-3-(4-trifluoromethylphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-amide;

N-{1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methylbenzenesulfonamide;

N-{1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-methoxy-N-methylbenzenesulfonamide;

N-[1-(3-biphenyl-4-yl-4-oxo-3,4-dihydroquinazolin-2-yl)ethyl]-4-tert-butyl-N-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(3'-methoxy-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(3'-chloro-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-{1-[3-(4'-methyl-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-benzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(2'-chloro-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4'-chloro-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(2'-methoxy-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4'-methoxy-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4'-(2-{1-[(4-*tert*-butylbenzenesulfonyl)methylamino]ethyl}-4-oxo-4*H*-quinazolin-3-yl)-biphenyl-4-carboxylic acid;

4-*tert*-butyl-*N*-{1-[3-(4'-cyano-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-thiophen-3-ylphenyl)-3,4-dihydroquinazolin-2-yl]ethyl} benzenesulfonamide;

4'-(2-{1-[(4-*tert*-butylbenzenesulfonyl)methylamino]ethyl}-4-oxo-4*H*-quinazolin-3-yl)-biphenyl-3-carboxylic acid methyl ester;

4'-(2-{1-[(4-*tert*-butylbenzenesulfonyl)methylamino]ethyl}-4-oxo-4*H*-quinazolin-3-yl)-biphenyl-4-carboxylic acid methyl ester;

4'-(2-{1-[(4-*tert*-butylbenzenesulfonyl)-methylamino]ethyl}-4-oxo-4*H*-quinazolin-3-yl)-biphenyl-3-carboxylic acid;

4-*tert*-butyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-pyrrolidin-1-ylphenyl)-3,4-dihydroquinazolin-2-yl]ethyl} benzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-piperidin-1-yl-phenyl)-3,4-dihydroquinazolin-2-yl]ethyl}benzenesulfonamide; and

4-*tert*-butyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-thiophen-2-ylphenyl)-3,4-dihydroquinazolin-2-yl]ethyl} benzenesulfonamide.

- 6. (Original) The compound of claim 3 wherein m is 0 or 1, n is 1, 2 or 3 and each R^{1a} is selected from alkyl.
- 7. (Currently amended) The compound of claim 6 selected from the group consisting of the following:

quinoline-8-sulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

naphthalene-1-sulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

naphthalene-2-sulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

2-naphthalen-1-yl-ethanesulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2,N-dimethylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-3,N-dimethylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,N-dimethylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methyl-C-phenyl-methanesulfonamide;

 $4-acetyl-N-\{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl\}-N-methylbenzenesulfonamide; \\$

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-3-trifluoromethylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methyl-4-trifluoromethoxy-benzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2,5,N-trimethylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-3,4-dimethoxy-N-methylbenzenesulfonamide;

N-[4-({1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-methyl-sulfamoyl)-phenyl]-acetamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2,4,6,N-tetramethylbenzenesulfonamide;

2-phenyl-ethenesulfonic acid {1-[3-(2,4-dimethylphenyl) 4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide; 2,2,5,6,8-pentamethyl-ehroman-7-sulfonic acid {1-[3-(2,4-dimethylphenyl) 4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-3,4-difluoro-N-methylbenzenesulfonamide;

3-chloro-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2,*N*-dimethylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-5-fluoro-2,N-dimethylbenzenesulfonamide;

3,5 dimethyl-isoxazole 4 sulfonic acid {1-[3-(2,4-dimethylphenyl) 4 oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-4-trifluoromethylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-3-fluoro-N-methylbenzenesulfonamide;

2,4,6-trichloro-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

3-chloro-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-fluoro-*N*-methylbenzenesulfonamide;

2-chloro-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

5-chloro-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2-methoxy-*N*-methylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2,5-dimethoxy-N-methylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2,3,4-trifluoro-*N*-methylbenzenesulfonamide;

 $3-chloro-N-\{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl\}-N-methylbenzenesulfonamide; \\$

4-cyano-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-butyl-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

 $\label{eq:N-local-equation} N-\{1-[3-(2,4-\text{dimethylphenyl})-4-\text{oxo-3},4-\text{dihydroquinazolin-2-yl}]\text{ethyl}\}-4-(1,1-\text{dimethyl-propyl})-N-\text{methylbenzenesulfonamide};$

4-butoxy-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-3-methoxy-*N*-methylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2-methoxy-4,N-dimethylbenzenesulfonamide;

4-chloro-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2,5,*N*-trimethylbenzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-methanesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-4-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-4-methoxy-benzenesulfonamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-methoxy-N-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-4-methoxy-N-methylbenzenesulfonamide;

4-chloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzenesulfonamide;

octane-1-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide; quinoline-8-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

naphthalene-1-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2,*N*-dimethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-3,*N*-dimethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-4,N-dimethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-N-methyl-3-trifluoromethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-N-methyl-4-trifluoromethoxy-benzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2,5,*N*-trimethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-3,4-dimethoxy-*N*-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2,4,6,*N*-tetramethylbenzenesulfonamide;

2,2,5,6,8-pentamethyl-chroman-7-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide; N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-3,4-difluoro-N-methylbenzenesulfonamide;

3-chloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2,*N*-dimethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-5-fluoro-2,N-dimethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-N-methyl-4-trifluoromethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-3-fluoro-*N*-methylbenzenesulfonamide;

2,4,6-trichloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzenesulfonamide;

3-chloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-4-fluoro-*N*-methylbenzenesulfonamide;

2-chloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzenesulfonamide;

5-chloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2-methoxy-*N*-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2,5-dimethoxy-*N*-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2,3,4-trifluoro-*N*-methylbenzenesulfonamide;

3-chloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

4-cyano-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzenesulfonamide;

4-butyl-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-4-(1,1-dimethyl-propyl)-N-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-4-isopropyl-N-methylbenzenesulfonamide;

4-butoxy-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-3-methoxy-N-methylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2-methoxy-4,N-dimethylbenzenesulfonamide;

4-chloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-2,5,*N*-trimethylbenzenesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-N-methyl-3,5-bis-trifluoromethylbenzenesulfonamide; and

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methyl-4-nitro-benzenesulfonamide;.

4-*tert*-butyl-*N*-{1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4,N-dimethylbenzenesulfonamide;

N-{1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-methoxy-N-methylbenzenesulfonamide;

N-{1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methylbenzenesulfonamide;

4-chloro-N- $\{1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl\}-<math>N$ -methylbenzenesulfonamide;

N-{1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-4-trifluoromethylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-*tert*-butyl-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-*tert*-butyl-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-*tert*-butyl-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-isopropyl-*N*-methyl-*N*-[1-(4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

biphenyl-4-sulfonic acid methyl-[1-(4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]amide;

4-*tert*-butyl-*N*-[1-(6-methoxy-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-[1-(6-hydroxy-4-oxo-3-p-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-*N*-methylbenzenesulfonamide;

N-[1-(6-bromo-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-4-*tert*-butyl-*N*-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid [1-(6-bromo-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]methylamide;

N-[1-(6-bromo-4-oxo-3-p-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-4-isopropyl-N-methylbenzenesulfonamide;

N-[1-(6-benzyloxy-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-4-*tert*-butyl-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-[1-(6-isobutoxy-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-*N*-methylbenzenesulfonamide;

N-[1-(6-butoxy-4-oxo-3-p-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-4-tert-butyl-N-methylbenzenesulfonamide;

(2-{1-[(4-*tert*-butylbenzenesulfonyl)methylamino]ethyl}-4-oxo-3-p-tolyl-3,4-dihydroquinazolin-6-yloxy)acetic acid ethyl ester;

4-*tert*-butyl-*N*-[1-(6-ethoxy-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-6-thiophen-3-yl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-6-phenyl-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-6-thiophen-2-yl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-{1-[3-(2'-methyl-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-benzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-{1-[3-(3'-methyl-biphenyl-4-yl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-benzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-6-o-tolyl-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-3,6-di-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-*tert*-butyl-*N*-{1-[6-(2-chlorophenyl)-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[6-(4-chlorophenyl)-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[6-(2-methoxyphenyl)-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[6-(3-methoxyphenyl)-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[6-(4-methoxyphenyl)-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

3-(2-{1-[(4-*tert*-butyl-benzenesulfonyl)methylamino]ethyl}-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-6-yl)benzoic acid methyl ester;

4-(2-{1-[(4-*tert*-butyl-benzenesulfonyl)methylamino]ethyl}-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-6-yl)-benzoic acid methyl ester;

3-(2-{1-[(4-*tert*-butylbenzenesulfonyl)methylamino]ethyl}-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-6-yl)benzoic acid;

4-*tert*-butyl-*N*-{1-[6-(4-cyanophenyl)-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(6-morpholin-4-yl-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide; and

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-6-m-tolyl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-6-pyrrolidin-1-yl-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-[2-{1-[(4-*tert*-butyl-benzenesulfonyl)methylamino]ethyl}-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-6-yl]benzoic acid;

4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-6-piperidin-1-yl-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide; and

4-*tert*-butyl-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide; <u>and</u>

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzenesulfonamide.

- 8. (Canceled)
- 9. (Canceled)

- 10. (Original) The compound of claim 3 wherein m is 0 or 1, n is 0 or 1 and each R la is selected from carboxy, dialkylamino, hydroxy, alkoxyalkoxy, alkoxycarbonylalkoxy, aralkoxy, and heterocyclylalkoxy.
- 11. (Original) The compound of claim 10 selected from the group consisting of the following:

4-(2-{1-[(4-*tert*-butyl-benzenesulfonyl)methylamino]ethyl}-4-oxo-4*H*-quinazolin-3-yl)-benzoic acid;

4-*tert*-butyl-*N*-{1-[3-(4-dimethylamino-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-dimethylamino-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-dimethylamino-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-isopropyl-*N*-methyl-*N*-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

biphenyl-4-sulfonic acid methyl-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]amide;

N-methyl-*N*-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-methoxy-*N*-methyl-*N*-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4,*N*-dimethyl-*N*-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-chloro-*N*-methyl-*N*-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

N-methyl-*N*-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]-4-trifluoromethylbenzenesulfonamide;

N-{1-[3-(4-benzyloxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-*tert*-butyl-*N*-methylbenzenesulfonamide;

N-{1-[3-(4-benzyloxy-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide;

biphenyl-4-sulfonic acid {1-[3-(4-benzyloxy-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(4-hydroxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-{1-[5-hydroxy-3-(4-hydroxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-(1-{4-oxo-3-[4-(2-piperidin-1-yl-ethoxy)phenyl]-3,4-dihydroquinazolin-2-yl}ethyl)benzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-(1-{3-[4-(2-morpholin-4-yl-ethoxy)phenyl]-4-oxo-3,4-dihydroquinazolin-2-yl}ethyl)benzenesulfonamide;

[4-(2-{1-[(4-*tert*-butylbenzenesulfonyl)methylamino]ethyl}-4-oxo-4*H*-quinazolin-3-yl)phenoxy]acetic acid ethyl ester; and

4-*tert*-butyl-*N*-(1-{3-[4-(2-methoxyethoxy)phenyl]-4-oxo-3,4-dihydroquinazolin-2-yl}ethyl)-*N*-methylbenzenesulfonamide.

- 12. (Canceled)
- 13. (Canceled)
- 14. (Canceled)
- 15. (Currently amended) The compound of Claim 2 wherein the compound of formula (II) is a compound of formula (V):

$$(R^3)_{m}$$
 $(R^{1a})_{n}$ (V)

wherein:

m is an integer from 0 to 4;

n is an integer from 0 to 5;

t is an integer from 0 to 5;

each R^{1a} is selected from the group consisting of alkyl, alkoxy, aralkoxy, halo, haloalkyl, haloalkoxy, amino, alkylamino, and dialkylamino;

R², R⁴ and R⁶ are each independently hydrogen or alkyl;

each R³ is independently selected from the group consisting of alkyl, alkoxy, and halo; and

each R^{5a} is independently selected from the group consisting of alkyl, alkoxy, alkoxycarbonyl, halo, and aryl;

or two adjacent R^{5a} groups form <u>phenyl</u>, 5-6 membered heteroaryl, O- $(CH_2)_y$ -O-, <u>-S- $(CH_2)_y$ -O-</u>, or -S- $(CH_2)_y$ -S-.; an aryl, heterocyclyl or heteroaryl

- 16. (Original) The compound of claim 15 wherein m is 0 or 1, n is 1 and each R^{1a} is alkoxy.
- 17. (Original) The compound of claim 16 selected from the group consisting of the following:

4-chloro-*N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide;

3-methoxy-*N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide;

4-methoxy-*N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide;

4-*tert*-butyl-*N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide;

N-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-terephthalamic acid methyl ester;

2,4-dichloro-*N*-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide;

benzo[1,3]dioxole-5-carboxylic acid [3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]amide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methyl-terephthalamic acid methyl ester;

2-methoxy-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

3-methoxy-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

4-methoxy-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-8-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-5-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[7-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-7-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[8-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-tert-butyl-N-{1-[8-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-

dihydroquinazolin-2-yl]ethyl}-N-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[5-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-tert-butyl-N-{1-[5-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-

dihydroquinazolin-2-yl]ethyl}-N-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[6-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[6-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

4-*tert*-butyl-*N*-{1-[3-(3-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(3-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(2-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(2-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[8-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[8-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide; and

4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide.

- 18. (Original) The compound of claim 15 wherein m is 0 or 1, n is 1, 2 or 3 and each R^{1a} is selected from alkyl.
- 19. (Original) The compound of claim 18 selected from the group consisting of the following:

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide; 4-tert-butyl-N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide;

benzo[1,3]dioxole-5-carboxylic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]amide;

2,4-dichloro-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]benzamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-terephthalamic acid methyl ester;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methyl-terephthalamic acid methyl ester;

benzo[1,3]dioxole-5-carboxylic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

4-*tert*-butyl-*N*-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-*N*-methylbenzamide;

4-*tert*-butyl-*N*-{1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(3,5-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(4-*tert*-butyl-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-*tert*-butyl-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl} methylamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-3-p-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-benzamide;

benzo[1,3]dioxole-5-carboxylic acid methyl-[1-(4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]amide;

benzo[1,3]dioxole-5-carboxylic acid [1-(6-bromo-4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]methylamide; and

N-[1-(6-bromo-4-oxo-3-p-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]-4-tert-butyl-N-methylbenzamide.

- 20. (Original) The compound of claim 15 wherein m is 0 or 1, n is 0 or 1 and each R^{1a} is independently selected from dialkylamino, aralkoxy, halo, haloalkyl and haloalkoxy.
- 21. (Original) The compound of claim 20 selected from the group consisting of the following:

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-dimethylamino-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-{1-[3-(4-dimethylamino-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-benzyloxy-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(4-benzyloxy-phenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-tert-butyl-N-methylbenzamide;

N-{1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-2-methoxy-N-methylbenzamide;

N-{1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-3-methoxy-N-methylbenzamide;

N-{1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-methoxy-N-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl} methylamide;

N-{1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methyl-terephthalamic acid methyl ester;

4-*tert*-butyl-*N*-{1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]benzamide;

benzo[1,3]dioxole-5-carboxylic acid methyl-[1-(4-oxo-3-phenyl-3,4-dihydroquinazolin-2-yl)ethyl]amide;

benzo[1,3]dioxole-5-carboxylic acid {1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(4-bromophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-tert-butyl-N-methylbenzamide;

4-*tert*-butyl-*N*-methyl-*N*-{1-[4-oxo-3-(4-trifluoromethoxyphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}benzamide; and

benzo[1,3]dioxole-5-carboxylic acid methyl-{1-[4-oxo-3-(4-trifluoromethoxyphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-amide;

4-*tert*-butyl-*N*-{1-[3-(4-fluorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzamide;

benzo[1,3]dioxole-5-carboxylic acid methyl-{1-[4-oxo-3-(4-trifluoromethylphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-amide; and 4-tert-butyl-N-methyl-N-{1-[4-oxo-3-(4-trifluoromethylphenyl)-3,4-dihydroquinazolin-2-yl]ethyl}-benzamide.

- 22. (Original) The compound of claim 1 wherein R¹ is hydrogen, optionally substituted alkyl, optionally substituted aryl, or optionally substituted aralkyl.
 - 23. (Canceled)
 - 24. (Canceled)
- 25. (Previously presented) The compound of claim 22 wherein R^1 is optionally substituted phenyl, and is selected with the proviso that it is not substituted at the 3- or 4-position with $-C(OH)(CF_3)_2$.
- 26. (Currently amended) The compound of claim 2 wherein any two R^{1a} or R^{3} groups, which substitute adjacent carbons on the ring, together form alkylene, alkenylene, $\underline{C_6}$ cycloalkylene, phenylene, alkylenedioxy, thioalkylenoxy, or alkylenedithioxy-O-($\underline{CH_2}$)_y-O-, $\underline{-S-(\underline{CH_2})_y-O-}$ or $\underline{-S-(\underline{CH_2})_y-S-}$, where y is 1 or 2.
 - 27. (Original) The compound of claim 2 wherein R^{1a} is not -C(OH)(CF₃)₂.
 - 28. (Original) The compound of claim 1 wherein R⁶ is hydrogen.
 - 29. (Canceled)

- 30. (Original) The compound of claim 1 wherein m is 1.
- 31. (Original) The compound of claim 2 wherein each R^{1a} is independently halo, pseudohalo, optionally substituted alkyl, optionally substituted alkoxy, optionally substituted aryl, optionally substituted dialkylamino, optionally substituted aralkoxy, hydroxy, optionally substituted heteroaryl, optionally substituted heterocyclyl or optionally substituted cycloalkyl.

32. (Canceled)

- 33. (Previously presented) The compound of claim 31 wherein each R^{1a} is independently chloro, fluoro, ethyl, methyl, methoxy, bromo, cyano, phenyl, *tert*-butyl, trifluoromethoxy, dimethylamino, trifluoromethyl, benzyloxy, hydroxy, 2-methylphenyl, 3-methylphenyl, 2-chlorophenyl, 3-chlorophenyl, 4-chlorophenyl, 2-methoxyphenyl, 3-methoxyphenyl, 4-methoxyphenyl, ethoxy, isopropoxy, butoxy, isobutoxy, 2-(*N*-morpholino)ethoxy, 2-methoxyethoxy, 4-cyanophenyl, 2-thienyl, 3-thienyl, 3-methoxycarbonylphenyl, 4-methoxycarbonylphenyl, 3-carboxyphenyl, *N*-pyrrolidinyl, or *N*-morpholinyl.
- 34. (Original) The compound of claim 1 wherein R^2 is hydrogen or optionally substituted alkyl, and R^6 is hydrogen.
 - 35. (Canceled)
- 36. (Previously presented) The compound of claim 34 wherein R² is hydrogen, methyl or ethyl.
- 37. (Original) The compound of claim 1 wherein each R³ is independently optionally substituted alkyl, halo, pseudohalo, optionally substituted alkoxy, hydroxy, optionally substituted aralkoxy, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted heterocyclyl, or optionally substituted cycloalkyl.

38. (Canceled)

- 39. (Previously presented) The compound of claim 37 wherein each R³ is independently methyl, chloro, methoxy, hydroxy, bromo, ethoxy, isopropoxy, isobutoxy, butoxy, benzyloxy, ethoxycarbonylmethoxy, phenyl, 2-thienyl, 3-thienyl, 2-methylphenyl, 3-methylphenyl, 4-methylpheny, 2-chlorophenyl, 3-chlorophenyl, 4-chlorophenyl, 3-methoxyphenyl, 4-methoxyphenyl, 4-carboxyphenyl, N-pyrrolidinyl, N-morpholinyl, 3-methoxycarbonylphenyl, 4-methoxycarbonylphenyl, 3-carboxyphenyl, 4-cyanophenyl, or piperidinyl.
- 40. (Original) The compound of claim 1 wherein one of R^4 and R^5 is -SO₂-(optionally substituted aryl).
- 41. (Original) The compound of claim 1 wherein one of R⁴ and R⁵ is -SO₂-(optionally substituted phenyl).
- 42. (Currently amended) The compound of claim 2 wherein the compound has formula (III):

$$(R^{3})_{m} = (R^{1a})_{n}$$

$$(R^{5a})_{t}$$

$$(R^{5a})_{t}$$

$$(R^{5a})_{t}$$

or a pharmaceutically acceptable derivative salt thereof, wherein:

t is an integer from 0 to 5;

R^{1a}, R², R³, R⁴, R⁶, n and m are as defined above;

each R^{5a} is independently optionally substituted alkyl, optionally substituted alkoxy, optionally substituted heteroaryl, optionally substituted aryl, optionally substituted heterocyclyl, halo, pseudohalo;

or any two R^{5a} substituents, which substitute adjacent atoms on the ring, together form phenyl, O-(CH₂)_v-O-, -S-(CH₂)_v-O-, or -S-(CH₂)_v-S-a optionally substituted cycloalkyl,

optionally substituted heterocyclyl, optionally substituted aryl, or optionally substituted heteroaryl ring having 5 or 6 members in the ring and where the heteroatoms, if present, are selected from O, S and optionally substituted N;

where R^{5a} , when substituted, is substituted with-one or more, in one embodiment one to five, in another embodiment one, two or three, substituents each independently selected from Q^2 - Q^4 , as defined above.

43. (Currently amended) The compound of claim 42 wherein any two R^{5a} groups, which substitute adjacent carbons on the ring, together form -N=C(R^{29})-C(R^{29})-C(R^{29})-or -C(R^{29})-C(R^{29})-C(R^{29})-,

where each R²⁹ is independently hydrogen, halo, pseudohalo, optionally substituted alkyl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aryl, optionally substituted aralkyl, or optionally substituted heteroaralkyl;

where R^{5a} and R^{29} , when substituted, are substituted with-one or more, in one embodiment one to five, in another embodiment one, two or three, substituents each independently selected from Q^2 Q^4 , as defined above.

44. (Canceled)

- 46. (Original) The compound of claim 42 wherein R⁴ is hydrogen, optionally substituted alkyl, optionally substituted aralkyl, or optionally substituted heteroaralkyl.

47. (Canceled)

- 48. (Previously presented) The compound of claim 46 wherein R⁴ is hydrogen, methyl, 2-methoxy-1-ethyl, propyl, isobutyl, butyl, pentyl, isopentyl, hexyl, benzyl, phenethyl or 2-thienylmethyl.
- 49. (Original) The compound of claim 42 wherein one R^{5a} group is 4-*tert*-butyl or 4-isopropyl.
- 50. (Currently amended) The compound of claim 2 wherein the compound has formula (IV):

$$(R^{3})_{m} = (R^{1a})_{n}$$

$$(IV)$$

$$R^{4} = (R^{5a})_{u}$$

$$R^{5b}$$

or a pharmaceutically acceptable derivative salt thereof, wherein R^{1a}, R², R³, R⁴, R⁶, R^{5a}, m and n are selected as above; wherein u is an integer from 0 to 4; and R^{5b} is *tert*-butyl or isopropyl.

- 51. (Canceled)
- 52. (Canceled)
- 53. (Original) The compound of claim 1 wherein one of R⁴ and R⁵ is -C(O)-(optionally substituted aryl).
- 54. (Original) The compound of claim 53 wherein one of R⁴ and R⁵ is -C(O)-(optionally substituted phenyl).

55. (Currently amended) The compound of claim 2 wherein the compound has formula (V):

$$(R^{3})_{m} = (R^{1a})_{n}$$

$$(V)$$

$$R^{4} = (R^{5a})_{t}$$

or a pharmaceutically acceptable derivative salt thereof, wherein R^{1a}, R², R³, R⁴, R⁶, R^{5a}, t, n and m are as defined above, wherein t is 0, 1, 2, 3, 4, or 5.

- 56. (Currently amended) The compound of claim 55 wherein each R^{1a} is independently halo, optionally substituted alkyl, or optionally substituted alkoxy, where the substituents, when present, are each independently selected from Q^{1} , as defined above.
 - 57. (Canceled)
- 58. (Previously presented) The compound of claim 56 wherein each R^{1a} is independently methoxy, methyl, chloro or fluoro.
- 59. (Currently amended) The compound of claim 55 wherein R^2 is hydrogen or optionally substituted alkyl, where the substituents, when present, are each independently selected from Q^1 , as defined above.
 - 60. (Canceled)
- 61. (Previously presented) The compound of claim 59 wherein R^2 is hydrogen or methyl.

- 62. (Currently amended) The compound of claim 55 wherein each R^3 is independently hydrogen or optionally substituted alkoxy, where the substituents, when present, are each independently selected from Q^1 , as defined above.
 - 63. (Canceled)
- 64. (Previously presented) The compound of claim 62 wherein each R³ is independently hydrogen or methoxy.
- 65. (Currently amended) The compound of claim 55 wherein R^4 is optionally substituted alkyl, where the substituents, when present, are each independently selected from Q^1 , as defined above.
 - 66. (Canceled)
- 67. (Previously presented) The compound of claim 65 wherein R⁴ is methyl or butyl.
 - 68. (Original) The compound of claim 55 wherein R⁶ is hydrogen.
- 69. (Currently amended) The compound of claim 55 wherein each R^{5a} is independently optionally substituted alkyl, where the substituents, when present, are each independently selected from Q^2Q^4 , as defined above.
 - 70. (Canceled)
- 71. (Previously presented) The compound of claim 69 wherein R^{5a} is *tert*-butyl.
- 72. (Currently amended) The compound of claim 1 wherein one of R^4 and R^5 is -C(O)-(optionally substituted alkyl), where the substituents, when present, are each independently selected from Q^1 , as defined above.
 - 73. (Canceled)

- 74. (Previously presented) The compound of claim 72 wherein one of R^4 and R^5 is -C(O)-octyl.
- 75. (Currently amended) The compound of claim 1 wherein one of R^4 and R^5 is $\underline{-C(O)-N(R^7)_2-C(O)-N(R^8)R^9}$, where R^8 and R^9 are each R^7 is independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclylalkyl;

or R^8 -and R^9 -together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene; where R^8 -and R^9 -are each R^7 is independently unsubstituted or substituted with one or more, in one embodiment one to five, in another embodiment one, two or three, substituents each independently selected from Q^1 , as defined above.

76. (Currently amended) The compound of claim 75 wherein \mathbb{R}^8 -and \mathbb{R}^9 -are each- \mathbb{R}^7 is independently selected from hydrogen, optionally substituted cycloalkyl, and optionally substituted aryl.

77. (Canceled)

78. (Currently amended) The compound of claim 76 wherein \mathbb{R}^8 -one \mathbb{R}^7 is hydrogen and \mathbb{R}^9 -the other \mathbb{R}^7 is cyclohexyl, 4-nitrophenyl, 2-methoxyphenyl, 3-cyanophenyl, 3,4-dichlorophenyl, 2,6-diisopropylphenyl, 2-methylphenyl, 2-trifluoromethylphenyl, 2-fluorophenyl, 3-fluorophenyl, 3-methylphenyl, 3-chlorophenyl, 2,6-dimethylphenyl or 3-trifluoromethylphenyl.

79. (Canceled)

80. (Previously presented) The compound of claim 1 wherein R⁴ and R⁵ together form -CH₂-C(H)(Me)-N(R³⁰)-CH₂-CH₂-, where R³⁰ is optionally substituted heteroarylcarbonyl, optionally substituted alkylcarbonyl, optionally substituted arylcarbonyl,

optionally substituted arylaminocarbonyl, or optionally substituted arylaminocarbonyl.

- 81. (Original) The compound of claim 80 wherein R³⁰ is 2-thienylcarbonyl, butyryl, 4-fluorobenzoyl, benzyloxyacetyl, diphenylacetyl, 4-nitrobenzoyl, 2,5-dichlorobenzenesulfonyl, *tert*-butylaminocarbonyl, phenylaminocarbonyl, 2,3-dichlorophenylaminocarbonyl, 4-*tert*-butylphenylsulfonyl or 3,4-methylenedioxybenzoyl.
 - 82. (Canceled)
- 83. (Previously presented) The compound of claim 1 wherein Q^2 is nitro, fluoro, benzyloxy or chloro; or two Q^2 groups, which substitute adjacent carbons, together form methylenedioxy.
- 84. (Previously presented) A method of treating, preventing, or ameliorating the symptoms of a disease or disorder that is modulated or otherwise affected by nuclear receptor activity or in which nuclear receptor activity is implicated, comprising administering to a subject in need thereof an effective amount of a compound of claim 1.
- 85. (Original) The method of claim 84, wherein the disease or disorder is selected from hypercholesterolemia, hyperlipoproteinemia, hypertriglyceridemia, lipodystrophy, hyperglycemia, diabetes mellitus, dyslipidemia, atherosclerosis, gallstone disease, acne vulgaris, acneiform skin conditions, diabetes, Parkinson's disease, cancer, Alzheimer's disease, inflammation, immunological disorders, lipid disorders, obesity, conditions characterized by a perturbed epidermal barrier function, hyperlipidemia, cholestasis, peripheral occlusive disease, ischemic stroke, conditions of disturbed differentiation or excess proliferation of the epidermis or mucous membrane, and cardiovascular disorders.
- 86. (Previously presented) A method of reducing cholesterol levels in a subject in need thereof, comprising administering an effective amount of a compound of claim 1.

- 87. (Previously presented) A method of treating, preventing, or ameliorating one or more symptoms of a disease or disorder which is affected by cholesterol, triglyceride, or bile acid levels, comprising administering to a subject in need thereof an effective amount of a compound of claim 1.
- 88. (Previously presented) A method of modulating nuclear receptor activity, comprising contacting the nuclear receptor with a compound of claim 1.
- 89. (Original) The method of claim 88, wherein the nuclear receptor is an orphan nuclear receptor.
- 90. (Original) The method of claim 88, wherein the nuclear receptor is farnesoid X receptor (FXR).
- 91. (Original) The method of claim 88, wherein the compound is a nuclear receptor agonist.
- 92. (Original) The method of claim 88, wherein the compound is a nuclear receptor antagonist.
- 93. (Previously presented) A method of modulating cholesterol metabolism, comprising administering an effective amount of a compound of claim 1.
- 94. (Previously presented) A method of treating, preventing or ameliorating one or more symptoms of hypocholesterolemia in a subject in need thereof, comprising administering an effective amount of a compound of claim 1.
- 95. (Previously presented) A method of increasing cholesterol efflux from cells of a subject, comprising administering an effective amount of a compound of claim 1.
- 96. (Previously presented) A method of increasing the expression of ATP-Binding Cassette (ABC1) in the cells of a subject, comprising administering an effective amount of a compound of claim 1.

- 97. (Previously presented) An *in vitro* method for altering nuclear receptor activity, comprising contacting the nuclear receptor with a compound of claim 1.
- 98. (Previously presented) The method of claim 84, wherein a second active agent selected from antihyperlipidemic agents, plasma HDL-raising agents, antihypercholesterolemic agents, cholesterol biosynthesis inhibitors (such as HMG CoA reductase inhibitors, such as lovastatin, simvastatin, pravastatin, fluvastatin, atorvastatin and rivastatin), acyl-coenzyme A:cholesterol acytransferase (ACAT) inhibitors, probucol, raloxifene, nicotinic acid, niacinamide, cholesterol absorption inhibitors, bile acid sequestrants (such as anion exchange resins, or quaternary amines (*e.g.*, cholestyramine or colestipol)), low density lipoprotein receptor inducers, clofibrate, fenofibrate, benzofibrate, cipofibrate, gemfibrizol, vitamin B₆, vitamin B₁₂, anti-oxidant vitamins, β-blockers, anti-diabetes agents, angiotensin II antagonists, angiotensin converting enzyme inhibitors, platelet aggregation inhibitors, fibrinogen receptor antagonists, aspirin and fibric acid derivatives; is administered simultaneously with, prior to, or after administration of the compound.
- 99. (Currently amended) A pharmaceutical composition, comprising, in a pharmaceutically acceptable carrier, a compound of formula (I):

$$(R^3)_{m} = \begin{pmatrix} 0 & R^1 & \\ N & R^6 & \\ N & R^2 & \\ N & R^5 & \\ N$$

wherein:

m is an integer from 0 to 4;

R¹ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroarylyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally

substituted heterocyclylalkyl, -OR⁷ or -N(R⁸)R⁹, with the proviso that R¹ is not 3- or 4- (1,1,1,3,3,3-hexafluoro-2-hydroxy-2-propyl)phenyl;

R², R⁴, R⁵ and R⁶ are selected from (a) and (b) as follows:

(a) R² and R⁶ are selected from (i) and (ii) as follows: (i) R² and R⁶ are each independently hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted heteroaryl, optionally substituted eycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, or optionally substituted heterocyclylalkyl; or (ii) R² and R⁶ together form optionally substituted alkylene or optionally substituted alkenylene; and

 R^4 and R^5 are selected from (i) and (ii) as follows: (i) R^4 and R^5 are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroarylyl, optionally substituted aralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-N(R^8)R^9$, $-OR^7$, $-S(O)_jR^{11}$ where j is 1 or 2, $-B(R^{22})_2$, $-P(R^{22})_2$, $-P(O)(R^{22})_2$ and $-C(E)R^{23}$, where E is selected from O, S and NR^7 ; or (ii) R^4 and R^5 together form optionally substituted alkylene, optionally substituted alkylene, optionally substituted alkylene; or

(b) R² and R⁵, or R² and R⁴, or R⁶ and R⁵, or R⁶ and R⁴, together form a 4, 5, 6 or 7 membered optionally substituted heterocyclyl group, or a 5 or 6 membered optionally substituted heteroaryl group; and the remainder of R², R⁴, R⁵ and R⁶ are each independently selected as in (i) above;

each R^3 is independently selected from the group consisting of halo, pseudohalo, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $N(R^{12})R^{13}$, $-OR^{14}$, $-C(E)R^{15}$ where E is O, S or NR^7 , and $-S(O)_yR^{16}$ where y is 0, 1 or 2;

or any two R³ groups, which substitute adjacent carbons on the ring, together form optionally substituted alkylene, optionally substituted alkenylene, optionally substituted alkylenedioxy, optionally substituted thioalkylenoxy, or optionally substituted alkylenedithioxy;

each R⁷ is independently selected from the group consisting of hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted recordly, optionally substituted eycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, and optionally substituted heterocyclylalkyl;

 R^8 and R^9 are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-S(O)_jR^{10}$ where j is 1 or 2, and $-C(M)R^{11}$, where M is selected from O and S;

or R⁸ and R⁹ together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene;

each R¹⁰ is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted eycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted eycloalkylalkyl, and optionally substituted heterocyclylalkyl;

each R^{11} is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-OR^{10}$ and $-N(R^7)_2$;

 R^{12} and R^{13} are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -C(M) R^{17} where M is O or S, and -S(O) $_iR^{18}$ where j is 1 or 2;

or R¹² and R¹³ together form optionally substituted alkylene, optionally substituted alkenylene, optionally substituted alkyleneoxyalkylene or optionally substituted alkyleneazaalkylene;

R¹⁴ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl or -C(M)R¹⁷ where M is O or S;

R¹⁵ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OH, -OR¹⁴ or -N(R¹²)R¹³;

R¹⁶ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OH, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁷ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OR¹⁹ or -N(R²⁰)R²¹;

 R^{18} is optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted aralkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted beterocyclylalkyl, optionally substituted heterocyclylalkyl, $-OR^{19}$ or $-N(R^{20})R^{21}$;

R¹⁹ is alkyl, alkenyl, alkynyl, cycloalkyl, heterocyclyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl or heteroaralkyl;

R²⁰ and R²¹ are each independently selected from hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, heterocyclyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl and heteroaralkyl,

or R^{20} and R^{21} together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene;

each R^{22} is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-OR^7$ and $-N(R^7)_2$;

 R^{23} is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-OR^{10}$, $-N(R^7)_2$, or $-N(R^7)N(R^7)_2$;

wherein each of the above R¹-R²³ groups, when substituted, are substituted with one or more substituents each independently selected from Q¹, where Q¹ is halo, pseudohalo, hydroxy, oxo, thia, nitrile, nitro, formyl, mercapto, carboxy, carboxyalkyl, alkyl, haloalkyl, polyhaloalkyl, aminoalkyl, diaminoalkyl, alkenyl containing 1 to 2 double bonds, alkynyl containing 1 to 2 triple bonds, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl, aralkenyl, aralkynyl, heteroarylalkyl, trialkylsilyl, dialkylarylsilyl,

alkyldiarylsilyl, triarylsilyl, alkylidene, arylalkylidene, alkylcarbonyl, cycloalkylcarbonyl, heterocyclylcarbonyl, arylcarbonyl, heteroarylcarbonyl, alkoxycarbonyl, alkoxycarbonylalkyl, aryloxycarbonyl, aryloxycarbonylalkyl, aralkoxycarbonyl, aralkoxycarbonylalkyl, arylcarbonylalkyl, aminocarbonyl, alkylaminocarbonyl, dialkylaminocarbonyl, arylaminocarbonyl, diarylaminocarbonyl, arylakylaminocarbonyl, alkoxy, aryloxy, heteroaryloxy, heteroaralkoxy, heterocyclyloxy, heterocyclylalkoxy, cycloalkoxy, perfluoroalkoxy, alkenyloxy, alkynyloxy, aralkoxy, alkylcarbonyloxy, arylcarbonyloxy, aralkylcarbonyloxy, alkoxycarbonyloxy, aryloxycarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, alkylarylaminocarbonyloxy, diarylaminocarbonyloxy, guanidino, isothioureido, amidino, alkylamidino, arylamidino, aminothiocarbonyl, alkylaminothiocarbonyl, arylaminothiocarbonyl, amino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, arylaminoalkyl, diarylaminoalkyl, alkylarylaminoalkyl, alkylamino, dialkylamino, haloalkylamino, arylamino, diarylamino, alkylarylamino, alkylcarbonylamino, alkoxycarbonylamino, aralkoxycarbonylamino, arylcarbonylamino, arylcarbonylaminoalkyl, aryloxycarbonylaminoalkyl, aryloxyarylcarbonylamino, aryloxycarbonylamino, alkylsulfonylamino, arylsulfonylamino, heteroarylsulfonylamino, heterocyclylsulfonylamino, heteroarylthio, azido, -N⁺(R²⁴)₃, -P(R²⁵)₂, $-P(O)(R^{25})_2$, $-OP(O)(R^{25})_2$, $-N(R^{24})C(O)R^{26}$, dialkylphosphonyl, alkylarylphosphonyl, diarylphosphonyl, hydroxyphosphonyl, alkylthio, arylthio, perfluoroalkylthio, carboxyalkylthio, thiocyano, isothiocyano, alkylsulfinyloxy, alkylsulfonyloxy, arylsulfinyloxy, arylsulfinyloxy, hydroxysulfonyloxy, alkoxysulfonyloxy, aminosulfonyloxy, alkylaminosulfonyloxy, dialkylaminosulfonyloxy, arylaminosulfonyloxy, diarylaminosulfonyloxy, alkylarylaminosulfonyloxy, alkylsulfinyl, alkylsulfonyl, arylsulfinyl, arylsulfonyl, hydroxysulfonyl, alkoxysulfonyl, aminosulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, arylaminosulfonyl, diarylaminosulfonyl or alkylarylaminosulfonyl; or two Q¹ groups, which substitute atoms in a 1,2 or 1,3 arrangement, together form alkylenedioxy (i.e., -O-(CH₂)_v-O-), thioalkylenoxy (i.e., -S-(CH₂)_v-O-) or alkylenedithioxy (i.e., -S-(CH₂)_v-S-) where y is 1 or 2; or two Q¹ groups, which substitute the same atom, together form alkylene;

each R²⁴ is independently selected from the group consisting of hydrogen, alkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, heterocyclyl and heterocyclylalkyl;

each R²⁵ is independently selected from the group consisting of hydroxy, alkoxy, aralkoxy, alkyl, heterocyclyl, aryl and -N(R²⁷)R²⁸,

 R^{26} is alkoxy, aralkoxy, alkyl, heteroaryl, heterocyclyl, aryl or -N(R^{27}) R^{28} ; R^{27} and R^{28} are each independently hydrogen, alkyl, aralkyl, aryl, heteroaryl, heteroaralkyl or heterocyclyl,

or R²⁷ and R²⁸ together form alkylene, azaalkylene, oxaalkylene or thiaalkylene; and each Q¹ is optionally substituted by one or more substituents selected from Q²; where each Q² is independently halo, pseudohalo, hydroxy, oxo, thia, nitrile, nitro, formyl, mercapto, carboxy, carboxyalkyl, alkyl, haloalkyl, polyhaloalkyl, aminoalkyl, diaminoalkyl, alkenyl containing 1 to 2 double bonds, alkynyl containing 1 to 2 triple bonds, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl, aralkenyl, aralkynyl, heteroarylalkyl, trialkylsilyl, dialkylarylsilyl, alkyldiarylsilyl, triarylsilyl, alkylidene, arylalkylidene, alkylcarbonyl, arylcarbonyl, heteroarylcarbonyl, alkoxycarbonyl, alkoxycarbonylalkyl, aryloxycarbonyl, aryloxycarbonylalkyl, aralkoxycarbonyl, aralkoxycarbonylalkyl, arylcarbonylalkyl, aminocarbonyl, alkylaminocarbonyl, dialkylaminocarbonyl, arylaminocarbonyl, diarylaminocarbonyl, arylalkylaminocarbonyl, alkoxy, aryloxy, heteroaryloxy, heteroaralkoxy, heterocyclyloxy, heterocyclylalkoxy, cycloalkoxy, perfluoroalkoxy, alkenyloxy, alkynyloxy, aralkoxy, alkylcarbonyloxy, arylcarbonyloxy, aralkylcarbonyloxy, alkoxycarbonyloxy, aryloxycarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, alkylarylaminocarbonyloxy, diarylaminocarbonyloxy, guanidino, isothioureido, amidino, alkylamidino, arylamidino, aminothiocarbonyl, alkylaminothiocarbonyl, arylaminothiocarbonyl, amino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, arylaminoalkyl, diarylaminoalkyl, alkylarylaminoalkyl, alkylamino, dialkylamino, haloalkylamino, arylamino, diarylamino, alkylarylamino, alkylcarbonylamino, alkoxycarbonylamino, aralkoxycarbonylamino, arylcarbonylamino, arylcarbonylaminoalkyl, aryloxycarbonylaminoalkyl, aryloxyarylcarbonylamino, aryloxycarbonylamino, alkylsulfonylamino, arylsulfonylamino,

heteroarylsulfonylamino, heterocyclylsulfonylamino, heteroarylthio, azido, $-N^+(R^{24})_3$, $-P(R^{25})_2$, $-P(O)(R^{25})_2$, $-OP(O)(R^{25})_2$, $-OP(O)(R^{25})_2$, $-N(R^{24})C(O)R^{26}$, dialkylphosphonyl, alkylarylphosphonyl, diarylphosphonyl, hydroxyphosphonyl, alkylthio, arylthio, perfluoroalkylthio, carboxyalkylthio, thiocyano, isothiocyano, alkylsulfinyloxy, alkylsulfonyloxy, arylsulfinyloxy, arylsulfonyloxy, hydroxysulfonyloxy, alkoxysulfonyloxy, aminosulfonyloxy, alkylaminosulfonyloxy, diarylaminosulfonyloxy, alkylsulfinyl, arylsulfinyl, arylsulfinyl, arylsulfonyl, hydroxysulfonyl, alkoxysulfonyl, alkylsulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, arylaminosulfonyl, diarylaminosulfonyl, arylsulfonyl, or two Q^2 groups, which substitute atoms in a 1,2 or 1,3 arrangement, together form alkylenedioxy (*i.e.*, -O-(CH₂)_y-O-), thioalkylenoxy (*i.e.*, -S-(CH₂)_y-O-) or alkylenedithioxy (*i.e.*, -S-(CH₂)_y-S-) where y is 1 or 2; or two Q^2 groups, which substitute the same atom, together form alkylene,

as a stereoisomer, racemate or mixture thereof; or as a pharmaceutically acceptable derivative salt thereof.

100. (Currently amended) A pharmaceutical composition, comprising, in a pharmaceutically acceptable carrier:

(i) a compound of formula (I):

wherein:

m is an integer from 0 to 4;

R¹ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted

aralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-OR^7$ or $-N(R^8)R^9$, with the proviso that R^1 is not 3- or 4- (1,1,1,3,3,3-hexafluoro-2-hydroxy-2-propyl)phenyl;

R², R⁴, R⁵ and R⁶ are selected from (a) and (b) as follows:

(a) R² and R⁶ are selected from (i) and (ii) as follows: (i) R² and R⁶ are each independently hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl; or (ii) R² and R⁶ together form optionally substituted alkylene or optionally substituted alkenylene; and

R⁴ and R⁵ are selected from (i) and (ii) as follows: (i) R⁴ and R⁵ are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -N(R⁸)R⁹), -OR⁷, -S(O)_jR¹¹ where j is 1 or 2, -B(R²²)₂, -P(R²²)₂, -P(O)(R²²)₂ and -C(E)R²³, where E is selected from O, S and NR⁷; or (ii) R⁴ and R⁵ together form optionally substituted alkylene, optionally substituted alkylene, optionally substituted alkyleneazaalkylene; or

(b) R² and R⁵, or R² and R⁴, or R⁶ and R⁵, or R⁶ and R⁴, together form a 4, 5, 6 or 7 membered optionally substituted heterocyclyl group, or a 5 or 6 membered optionally substituted heteroaryl group; and the remainder of R², R⁴, R⁵ and R⁶ are each independently selected as in (i) above;

each R³ is independently selected from the group consisting of halo, pseudohalo, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted

heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, - $N(R^{12})R^{13}$, - OR^{14} , - $C(E)R^{15}$ where E is O, S or NR^7 , and - $S(O)_yR^{16}$ where y is 0, 1 or 2;

or any two R³ groups, which substitute adjacent carbons on the ring, together form optionally substituted alkylene, optionally substituted alkenylene, optionally substituted alkylenedioxy, optionally substituted thioalkylenoxy, or optionally substituted alkylenedithioxy;

each R⁷ is independently selected from the group consisting of hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted recordly, optionally substituted eycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, and optionally substituted heterocyclylalkyl;

 R^8 and R^9 are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-S(O)_jR^{10}$ where j is 1 or 2, and $-C(M)R^{11}$, where M is selected from O and S;

or R⁸ and R⁹ together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene;

each R¹⁰ is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted eycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted eycloalkylalkyl, and optionally substituted heterocyclylalkyl;

each R¹¹ is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted recordly, optionally substituted heteroaryl, optionally substituted heteroaralkyl,

optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-OR^{10}$ and $-N(R^7)_2$;

 R^{12} and R^{13} are each independently selected from hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heteroaralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -C(M) R^{17} where M is O or S, and -S(O) $_iR^{18}$ where j is 1 or 2;

or R¹² and R¹³ together form optionally substituted alkylene, optionally substituted alkenylene, optionally substituted alkyleneoxyalkylene or optionally substituted alkyleneazaalkylene;

 R^{14} is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl or $-C(M)R^{17}$ where M is O or S;

R¹⁵ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OH, -OR¹⁴ or -N(R¹²)R¹³;

R¹⁶ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OH, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁷ is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted

aralkyl, optionally substituted heteroaralkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁸ is optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted aralkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted beterocyclylalkyl, optionally substituted heterocyclylalkyl, -OR¹⁹ or -N(R²⁰)R²¹;

R¹⁹ is alkyl, alkenyl, alkynyl, cycloalkyl, heterocyclyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl or heteroaralkyl;

R²⁰ and R²¹ are each independently selected from hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, heterocyclyl, cycloalkylalkyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl and heteroaralkyl,

or R^{20} and R^{21} together form alkylene, alkenylene, alkyleneoxyalkylene or alkyleneazaalkylene;

each R^{22} is independently selected from the group consisting of optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, $-OR^7$ and $-N(R^7)_2$;

 R^{23} is hydrogen, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted aryl, optionally substituted heteroaryl, optionally substituted cycloalkyl, optionally substituted heterocyclyl, optionally substituted aralkyl, optionally substituted cycloalkylalkyl, optionally substituted cycloalkylalkyl, optionally substituted heterocyclylalkyl, $-OR^{10}$, $-N(R^7)_2$, or $-N(R^7)N(R^7)_2$;

wherein each of the above R¹-R²³ groups, when substituted, are substituted with one or more substituents each independently selected from Q¹, where Q¹ is halo, pseudohalo, hydroxy, oxo, thia, nitrile, nitro, formyl, mercapto, carboxy, carboxyalkyl, alkyl, haloalkyl, polyhaloalkyl, aminoalkyl, diaminoalkyl, alkenyl containing 1 to 2 double bonds, alkynyl

containing 1 to 2 triple bonds, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl, aralkenyl, aralkynyl, heteroarylalkyl, trialkylsilyl, dialkylarylsilyl, alkyldiarylsilyl, triarylsilyl, alkylidene, arylalkylidene, alkylcarbonyl, cycloalkylcarbonyl, heterocyclylcarbonyl, arylcarbonyl, heteroarylcarbonyl, alkoxycarbonyl, alkoxycarbonylalkyl, aryloxycarbonyl, aryloxycarbonylalkyl, aralkoxycarbonyl, aralkoxycarbonylalkyl, arylcarbonylalkyl, aminocarbonyl, alkylaminocarbonyl, dialkylaminocarbonyl, arylaminocarbonyl, diarylaminocarbonyl, arylalkylaminocarbonyl, alkoxy, aryloxy, heteroaryloxy, heteroaralkoxy, heterocyclyloxy, heterocyclylalkoxy, cycloalkoxy, perfluoroalkoxy, alkenyloxy, alkynyloxy, aralkoxy, alkylcarbonyloxy, arylcarbonyloxy, aralkylcarbonyloxy, alkoxycarbonyloxy, aryloxycarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, alkylarylaminocarbonyloxy, diarylaminocarbonyloxy, guanidino, isothioureido, amidino, alkylamidino, arylamidino, aminothiocarbonyl, alkylaminothiocarbonyl, arylaminothiocarbonyl, amino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, arylaminoalkyl, diarylaminoalkyl, alkylarylaminoalkyl, alkylamino, dialkylamino, haloalkylamino, arylamino, diarylamino, alkylarylamino, alkylcarbonylamino, alkoxycarbonylamino, aralkoxycarbonylamino, arylcarbonylamino, arylcarbonylaminoalkyl, aryloxycarbonylaminoalkyl, aryloxyarylcarbonylamino, aryloxycarbonylamino, alkylsulfonylamino, arylsulfonylamino, heteroarylsulfonylamino, heterocyclylsulfonylamino, heteroarylthio, azido, -N⁺(R²⁴)₃, -P(R²⁵)₂, $-P(O)(R^{25})_2, -OP(O)(R^{25})_2, -N(R^{24})C(O)R^{26}, \\ dialkylphosphonyl, \\ alkylarylphosphonyl, \\$ diarylphosphonyl, hydroxyphosphonyl, alkylthio, arylthio, perfluoroalkylthio, carboxyalkylthio, thiocyano, isothiocyano, alkylsulfinyloxy, alkylsulfonyloxy, arylsulfinyloxy, arylsulfinyloxy, hydroxysulfonyloxy, alkoxysulfonyloxy, aminosulfonyloxy, alkylaminosulfonyloxy, dialkylaminosulfonyloxy, arylaminosulfonyloxy, diarylaminosulfonyloxy, alkylarylaminosulfonyloxy, alkylsulfinyl, alkylsulfonyl, arylsulfinyl, arylsulfonyl, hydroxysulfonyl, alkoxysulfonyl, aminosulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, arylaminosulfonyl, diarylaminosulfonyl or alkylarylaminosulfonyl; or two O¹ groups, which substitute atoms in a 1,2 or 1,3 arrangement, together form alkylenedioxy (i.e., -O-(CH₂)_v-O-),

thioalkylenoxy (*i.e.*, -S-(CH₂)_y-O-) or alkylenedithioxy (*i.e.*, -S-(CH₂)_y-S-) where y is 1 or 2; or two Q^1 groups, which substitute the same atom, together form alkylene;

each R²⁴ is independently selected from the group consisting of hydrogen, alkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, heterocyclyl and heterocyclylalkyl;

each R²⁵ is independently selected from the group consisting of hydroxy, alkoxy, aralkoxy, alkyl, heterocyclyl, aryl and -N(R²⁷)R²⁸,

 R^{26} is alkoxy, aralkoxy, alkyl, heteroaryl, heterocyclyl, aryl or -N(R^{27}) R^{28} ; R^{27} and R^{28} are each independently hydrogen, alkyl, aralkyl, aryl, heteroaryl, heteroaralkyl or heterocyclyl,

or R²⁷ and R²⁸ together form alkylene, azaalkylene, oxaalkylene or thiaalkylene; and each Q¹ is optionally substituted by one or more substituents selected from Q²; where each Q² is independently halo, pseudohalo, hydroxy, oxo, thia, nitrile, nitro, formyl, mercapto, carboxy, carboxyalkyl, alkyl, haloalkyl, polyhaloalkyl, aminoalkyl, diaminoalkyl, alkenyl containing 1 to 2 double bonds, alkynyl containing 1 to 2 triple bonds, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, heteroaryl, aralkyl, aralkynyl, aralkynyl, heteroarylalkyl, trialkylsilyl, dialkylarylsilyl, alkyldiarylsilyl, triarylsilyl, alkylidene, arylalkylidene, alkylcarbonyl, arylcarbonyl, heteroarylcarbonyl, alkoxycarbonyl, alkoxycarbonylalkyl, aryloxycarbonyl, aryloxycarbonylalkyl, aralkoxycarbonyl, aralkoxycarbonylalkyl, arylcarbonylalkyl, aminocarbonyl, alkylaminocarbonyl, dialkylaminocarbonyl, arylaminocarbonyl, diarylaminocarbonyl, arylalkylaminocarbonyl, alkoxy, aryloxy, heteroaryloxy, heteroaralkoxy, heterocyclyloxy, heterocyclylalkoxy, cycloalkoxy, perfluoroalkoxy, alkenyloxy, alkynyloxy, aralkoxy, alkylcarbonyloxy, arylcarbonyloxy, aralkylcarbonyloxy, alkoxycarbonyloxy, aryloxycarbonyloxy, aralkoxycarbonyloxy, aminocarbonyloxy, alkylaminocarbonyloxy, dialkylaminocarbonyloxy, alkylarylaminocarbonyloxy, diarylaminocarbonyloxy, guanidino, isothioureido, amidino, alkylamidino, arylamidino, aminothiocarbonyl, alkylaminothiocarbonyl, arylaminothiocarbonyl, amino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, arylaminoalkyl, diarylaminoalkyl, alkylarylamino, dialkylamino, haloalkylamino, arylamino, diarylamino, alkylarylamino, alkylcarbonylamino, alkoxycarbonylamino, aralkoxycarbonylamino,

arylcarbonylamino, arylcarbonylaminoalkyl, aryloxycarbonylaminoalkyl, aryloxyarylcarbonylamino, arylcarbonylamino, arylcarbonylamino, arylcarbonylamino, arylcarbonylamino, arylcarbonylamino, heteroarylcarbonylamino, heteroarylcarbonylamino, heteroarylcarbonylamino, heteroarylcarbonylamino, heteroarylcarbonylamino, heteroarylcarbonylamino, heteroarylcarbonylamino, heteroarylcarbonylcarbonylamino, heteroarylcarbonylcarbo

where R^{24} , R^{25} , R^{26} , R^{27} and R^{28} are as defined above;

as a stereoisomer, racemate or mixture thereof; or as a pharmaceutically acceptable derivative salt thereof; and

(ii) one or more of a second active agent selected from antihyperlipidemic agents, plasma HDL-raising agents, antihypercholesterolemic agents, cholesterol biosynthesis inhibitors (such as HMG CoA reductase inhibitors, such as lovastatin, simvastatin, pravastatin, fluvastatin, atorvastatin and rivastatin), acyl-coenzyme A:cholesterol acytransferase (ACAT) inhibitors, probucol, raloxifene, nicotinic acid, niacinamide, cholesterol absorption inhibitors, bile acid sequestrants (such as anion exchange resins, or quaternary amines (e.g., cholestyramine or colestipol)), low density lipoprotein receptor inducers, clofibrate, fenofibrate, benzofibrate, cipofibrate, gemfibrizol, vitamin B_6 , vitamin B_{12} , anti-oxidant vitamins, β -blockers, LXR \square or \square agonists or antagonists, anti-diabetes agents, angiotensin II antagonists, angiotensin converting enzyme inhibitors, platelet aggregation inhibitors, fibrinogen receptor antagonists, aspirin and fibric acid derivatives.

101. (Previously presented) A compound of Claim 1 selected from the group consisting of the following:

2-naphthalen-1-yl-ethanesulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-C-phenyl-methanesulfonamide;

2-phenyl-ethenesulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

thiophene-2-sulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

C-(7,7-dimethyl-2-oxo-bicyclo[2.2.1]hept-1-yl)-N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methyl-methanesulfonamide;

3,5-dimethyl-isoxazole-4-sulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl} methylamide;

N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-N-methylmethanesulfonamide;

octane-1-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

2-naphthalen-1-yl-ethanesulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyllmethylamide;

2-phenyl-ethenesulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

thiophene-2-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

C-(7,7-dimethyl-2-oxo-bicyclo[2.2.1]hept-1-yl)-N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-N-methyl-methanesulfonamide;

N-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-N-methylmethanesulfonamide;

butane-1-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

3-chloropropane-1-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide;

[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]-carbamic acid benzyl ester;

nonanoic acid {1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

nonanoic acid [3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]amide;

nonanoic acid {1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

nonanoic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(3-methyl-4-oxo-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-isopropyl-*N*-methyl-*N*-[1-(3-methyl-4-oxo-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

biphenyl-4-sulfonic acid methyl-[1-(3-methyl-4-oxo-3,4-dihydroquinazolin-2-yl)ethyl]amide;

4-methoxy-*N*-methyl-*N*-[1-(3-methyl-4-oxo-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-chloro-*N*-methyl-*N*-[1-(3-methyl-4-oxo-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;

4-*tert*-butyl-*N*-methyl-*N*-[1-(3-methyl-4-oxo-3,4-dihydroquinazolin-2-yl)ethyl]benzamide;

3-(4-methoxyphenyl)-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one;

3-(3,5-dimethylphenyl)-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one;

2-(1-methylaminoethyl)-3-phenyl-3*H*-quinazolin-4-one;

- 3-(4-bromophenyl)-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one;
- 3-(4-methoxyphenyl)-7-methyl-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one;
- 8-methoxy-3-(4-methoxyphenyl)-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one;
- 5-methoxy-3-(4-methoxyphenyl)-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one;
- 6-methoxy-3-(4-methoxyphenyl)-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one;
- 3-(4-dimethylamino-phenyl)-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one;
- 3-(4-fluorophenyl)-2-(1-methylaminoethyl)-3*H*-quinazolin-4-one
- 4-*tert*-butyl-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;
- 4-*tert*-butyl-*N*-{1-[6-methoxy-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;
- 4-*tert*-butyl-*N*-{1-[3-(4-chlorophenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;
- 4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;
- 4-*tert*-butyl-*N*-methyl-*N*-[1-(4-oxo-3-*p*-tolyl-3,4-dihydroquinazolin-2-yl)ethyl]benzenesulfonamide;
- 4-*tert*-butyl-*N*-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;
- 4-*tert*-butyl-*N*-{1-[3-(4-methoxyphenyl)-6-methyl-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;
- 4-*tert*-butyl-*N*-{1-[6-chloro-3-(4-methoxyphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-*N*-methylbenzenesulfonamide;
- N-{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}-4-isopropyl-N-methylbenzenesulfonamide; and
- biphenyl-4-sulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide.
 - 102. (Canceled)

- 103. (Previously presented) The compound of claim 50 having one or more features selected from the group consisting of the following:
- a) R^{1a} is selected from the group consisting of alkyl, hydroxy, alkoxy, alkoxy, and aralkoxy;
 - b) R⁴ is hydrogen or alkyl;
- c) R³ is independently selected from the group consisting of hydrogen, alkyl, alkoxy, and halo; and
- d) R^{5a} is independently selected from the group consisting of alkyl, alkoxy, halo, and alkylcarbonyl.
- 104. (Previously presented) The compound of claim 103 having one or more features selected from the group consisting of the following:
 - a) R^{1a} is alkoxy;
 - b) R⁴ is alkyl;
- c) R³ is selected from the group consisting of hydrogen, halo, alkoxy, and alkyl; and
 - d) R^{5a} is selected from the group consisting of *tert*-butyl and isopropyl.
- 105. (New) The compound of claim 2 selected from the group consisting of the following:

2-phenyl-ethenesulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

3,5-dimethyl-isoxazole-4-sulfonic acid {1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl}methylamide;

 $N-\{1-[3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-yl]ethyl\}-3-methoxy-<math>N$ -methylbenzenesulfonamide; and

octane-1-sulfonic acid [3-(2,4-dimethylphenyl)-4-oxo-3,4-dihydroquinazolin-2-ylmethyl]methylamide.